

Listing of Claims:

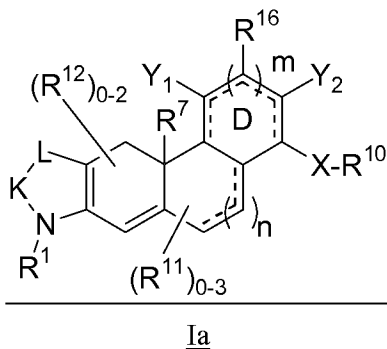
The listing of the claims which follows replaces any and all prior versions and/or listings of the claims in the application.

1-4. (Canceled)

5-10. (Previously Canceled)

11-18. (Canceled)

19. (Currently Amended) The A compound ~~according to Claim 18~~ represented by
Formula Ia:



or a pharmaceutically acceptable salt or hydrate thereof, wherein

R¹⁰ is selected from the group consisting of:

- (1) phenyl,
- (2) benzyl, and
- (3) HET, wherein HET is a 5-membered aromatic or non-aromatic monocyclic ring having 1-3 heteroatoms selected from O, S and N,

wherein groups (1) to (3) above are optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (a) halo,
- (b) C₁₋₄alkyl, optionally substituted with hydroxy or 1 to 3 halo groups,
- (c) C₁₋₄alkoxy, optionally substituted with 1 to 3 halo groups,
- (d) NH₂,
- (e) hydroxy, and
- (e) phenyl or benzyl;[. .]

n and m are each independently 0, 1 or 2;

K is selected from NR³ or C(R³)(R⁴);

L is selected from NR⁵ or C(R⁵)(R⁶);

X is a bond, -C(O), -N(R¹⁴)-, -N(R¹⁴)-C(O)-, -C(O)-N(R¹⁴)-, -N(R¹⁴)-S(O)_k-, -N(R¹⁴)-C(O)-NH- or -S(O)_k-N(R¹⁴);

k is 0, 1 or 2;

R¹ is selected from the group consisting of:

- (1) C₁₋₆alkyl,
- (2) C₂₋₆alkenyl,
- (3) C₂₋₆alkynyl,
- (4) C₃₋₆cycloalkyl,
- (5) C₁₋₆alkoxy,
- (6) C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2,
- (7) aryl,
- (8) aryl C₁₋₆alkyl,
- (9) HET,
- (10) -C₁₋₆alkyl-HET,
- (11) aryloxy,
- (12) aroyloxy,
- (13) aryl C₂₋₆alkenyl,
- (14) aryl C₂₋₆alkynyl,
- (15) hydrogen,
- (16) hydroxyl and
- (17) cyano

wherein items (1) to (6) above and the alkyl portions of items (8) and (10) above and the alkenyl portion of item (13) above and the alkynyl portion of item (14) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, oxo, OR¹³, N(R¹⁴)₂, C₃₋₆cycloalkyl and C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2, and

wherein the aryl is optionally substituted from one up to the maximum number of substitutable positions with halo; the aryl is items (7), (9), (11) and (12) above and aryl portion of items (8), (13) and (14) above and the HET portion of item (10) above are optionally substituted

from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of:

- (a) halo,
- (b) OR¹³,
- (c) N(R¹⁴)₂,
- (d) C₁₋₆alkyl,
- (e) C₂₋₆alkenyl,
- (f) C₂₋₆akynyl,
- (g) C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2,
- (h) aryl,
- (i) aryl-S(O)_k-, wherein k is 0, 1 or 2,
- (j) HET,
- (k) aryl C₁₋₆alkyl,
- (l) aroyl,
- (m) aryloxy,
- (n) aryl C₁₋₆alkoxy,
- (o) CN and
- (p) C₃₋₆cycloalkyl,

wherein items (d) to (g) and (p) above and the alkyl portions of item (k) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³ and N(R¹⁴)₂, and

wherein items (h), (i), (j), (l) and (m) above and the aryl portions of items (k) and (n) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³ and C₁₋₄alkyl,

R³, R⁴, R⁵ and R⁶ are each independently selected from the group consisting of:

- (1) hydrogen,
- (2) halo,
- (3) C₁₋₆alkyl,
- (4) C₂₋₆alkenyl,
- (5) C₂₋₆akynyl,
- (6) C₃₋₆cycloalkyl,
- (7) C₁₋₆alkoxy,
- (8) C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2,
- (9) aryl,
- (10) aryl C₁₋₆alkyl,
- (11) HET and

(12) -C₁₋₆alkyl-HET,

wherein items (3) to (8) above and the alkyl portions of items (10) and (12) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³, N(R¹⁴)₂ and C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2; and

wherein items (9) and (11) and the aryl portion of items (10) and the HET portion of item (12) are optionally substituted from one up to the maximum number of substituable positions with a substituent independently selected from the group consisting of:

- (a) halo,
- (b) OR¹³,
- (c) N(R¹⁴)₂,
- (d) C₁₋₆alkyl,
- (e) C₂₋₆alkenyl,
- (f) C₂₋₆alkynyl and
- (g) C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2,

wherein items (d) to (g) above are optionally substituted with from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³ and N(R¹⁴)₂,

or R¹ and R³ or R³ and R⁵ may be joined together to form a double bond;

R⁷ is selected from the group consisting of:

- (1) hydrogen,
- (2) OR¹³,
- (3) C₁₋₄alkyl,
- (4) aryl and
- (5) aryl C₁₋₄alkyl,

wherein item (3) above and the alkyl portion of item (5) above are optionally substituted with from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³ and N(R¹⁴)₂, and

wherein item (4) above and the aryl portion of item (5) above are optionally substituted with from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of:

- (a) halo,
- (b) OR¹³,
- (c) N(R¹⁴)₂,

- (d) C₁₋₆alkyl,
- (e) C₂₋₆alkenyl and
- (f) C₂₋₆alkynyl,

wherein items (d) to (f) above are optionally substituted with from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³ and N(R¹⁴)₂;

Y₁ is selected from the group consisting of:

- (1) hydrogen,
- (2) -O-R⁹,
- (3) -S(O)_k-R⁹, wherein k is 0, 1 or 2,
- (4) -C-W-R⁹, wherein W is O or S(O)_k,
- (5) -N(R¹⁵)₂,
- (6) -S(O)_k-N(R¹⁵)₂,
- (7) -N(R¹⁵)-S(O)_k-N(R¹⁵)₂,
- (8) NO₂,
- (9) -C(O)-R¹⁵,
- (10) -C(O)O-R¹⁵,
- (11) -CN,
- (12) halo,
- (13) -O-S(O)_k-R¹⁵ and
- (14) C₁₋₄alkyl, optionally substituted with from 1 to 6 halo groups,

Y₂ is CF₃;

R⁹ is selected from the group consisting of: hydrogen, C₁₋₁₂alkyl and aryl, wherein C₁₋₁₂alkyl and aryl are optionally substituted from one up to the maximum number of substituents with halo;

each R¹¹, R¹² and R¹⁶ is independently selected from the group consisting of:

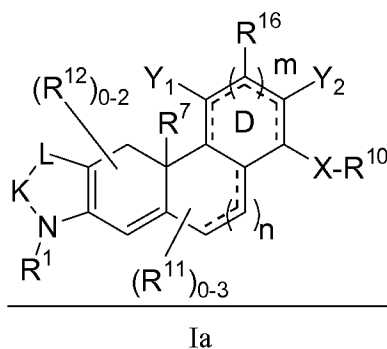
- (1) hydrogen,
- (2) halo,
- (3) C₁₋₆alkyl,
- (4) C₂₋₆alkenyl,
- (5) C₁₋₆alkoxy and
- (6) hydroxy,

wherein items (3) to (5) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR₁₂, N(R¹³)₂ and C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2,

each R¹³ and R¹⁴ is independently selected from the group consisting of hydrogen and C₁₋₄alkyl, optionally substituted from one up to the maximum number of substitutable positions with halo; and

each R¹⁵ is independently selected from the group consisting of: hydrogen, C₁₋₆alkyl, aryl and C₁₋₁₂alkoxycarbonyl, wherein said C₁₋₆alkyl and C₁₋₁₂alkoxycarbonyl are optionally substituted from one up to the maximum number of substitutable positions with halo and said aryl is optionally substituted from one up to the maximum number of substitutable positions with halo and C₁₋₄alkyl, optionally substituted with 1-3 halo groups.

20. (Currently Amended) ~~The~~ A compound ~~according to Claim 3~~ represented by Formula Ia:



or a pharmaceutically acceptable salt or hydrate thereof, wherein

Y₂ is hydrogen, X is a bond, ~~and~~ R¹⁰ is HET, wherein HET is a 5-membered aromatic or non-aromatic monocyclic ring containing 1-3 heteroatoms selected from O, S and N;[[.]]

n and m are each independently 0, 1 or 2;

K is selected from NR³ or C(R³)(R⁴);

L is selected from NR⁵ or C(R⁵)(R⁶);

R¹ is selected from the group consisting of:

(1) C₁₋₆alkyl,

- (2) C₂₋₆alkenyl,
- (3) C₂₋₆alkynyl,
- (4) C₃₋₆cycloalkyl,
- (5) C₁₋₆alkoxy,
- (6) C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2,
- (7) aryl,
- (8) aryl C₁₋₆alkyl,
- (9) HET,
- (10) -C₁₋₆alkyl-HET,
- (11) aryloxy,
- (12) aryloxy,
- (13) aryl C₂₋₆alkenyl,
- (14) aryl C₂₋₆alkynyl,
- (15) hydrogen,
- (16) hydroxyl and
- (17) cyano

wherein items (1) to (6) above and the alkyl portions of items (8) and (10) above and the alkenyl portion of item (13) above and the alkynyl portion of item (14) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, oxo, OR¹³, N(R¹⁴)₂, C₃₋₆cycloalkyl and C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2, and

wherein the aryl is optionally substituted from one up to the maximum number of substitutable positions with halo; the aryl is items (7), (9), (11) and (12) above and aryl portion of items (8), (13) and (14) above and the HET portion of item (10) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of:

- (a) halo,
- (b) OR¹³,
- (c) N(R¹⁴)₂,
- (d) C₁₋₆alkyl,
- (e) C₂₋₆alkenyl,
- (f) C₂₋₆alkynyl,
- (g) C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2,
- (h) aryl,
- (i) aryl-S(O)_k-, wherein k is 0, 1 or 2,
- (j) HET,

- (k) aryl C₁₋₆alkyl,
- (l) aroyl,
- (m) aryloxy,
- (n) aryl C₁₋₆alkoxy,
- (o) CN and
- (p) C₃₋₆cycloalkyl,

wherein items (d) to (g) and (p) above and the alkyl portions of item (k) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³ and N(R¹⁴)₂, and

wherein items (h), (i), (j), (l) and (m) above and the aryl portions of items (k) and (n) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³ and C₁₋₄alkyl,

R³, R⁴, R⁵ and R⁶ are each independently selected from the group consisting of:

- (1) hydrogen,
- (2) halo,
- (3) C₁₋₆alkyl,
- (4) C₂₋₆alkenyl,
- (5) C₂₋₆alkynyl,
- (6) C₃₋₆cycloalkyl,
- (7) C₁₋₆alkoxy,
- (8) C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2,
- (9) aryl,
- (10) aryl C₁₋₆alkyl,
- (11) HET and
- (12) -C₁₋₆alkyl-HET,

wherein items (3) to (8) above and the alkyl portions of items (10) and (12) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³, N(R¹⁴)₂ and C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2; and

wherein items (9) and (11) and the aryl portion of items (10) and the HET portion of item (12) are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of:

- (a) halo,
- (b) OR¹³,
- (c) N(R¹⁴)₂,

- (d) C₁₋₆alkyl,
- (e) C₂₋₆alkenyl,
- (f) C₂₋₆alkynyl and
- (g) C₁₋₆alkyl-S(O)_k, wherein k is 0, 1 or 2,

wherein items (d) to (g) above are optionally substituted with from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³ and N(R¹⁴)₂,

or R¹ and R³ or R³ and R⁵ may be joined together to form a double bond;

R⁷ is selected from the group consisting of:

- (1) hydrogen,
- (2) OR¹³,
- (3) C₁₋₄alkyl,
- (4) aryl and
- (5) aryl C₁₋₄alkyl,

wherein item (3) above and the alkyl portion of item (5) above are optionally substituted with from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³ and N(R¹⁴)₂, and

wherein item (4) above and the aryl portion of item (5) above are optionally substituted with from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of:

- (a) halo,
- (b) OR¹³,
- (c) N(R¹⁴)₂,
- (d) C₁₋₆alkyl,
- (e) C₂₋₆alkenyl and
- (f) C₂₋₆alkynyl,

wherein items (d) to (f) above are optionally substituted with from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹³ and N(R¹⁴)₂;

Y₁ is selected from the group consisting of:

- (1) hydrogen,
- (2) -O-R⁹,
- (3) -S(O)_k-R⁹, wherein k is 0, 1 or 2,
- (4) -C-W-R⁹, wherein W is O or S(O)_k,

- (5) -N(R¹⁵)₂,
- (6) -S(O)_k-N(R¹⁵)₂,
- (7) -N(R¹⁵)-S(O)_k-N(R¹⁵)₂,
- (8) NO₂,
- (9) -C(O)-R¹⁵,
- (10) -C(O)O-R¹⁵,
- (11) -CN,
- (12) halo,
- (13) -O-S(O)_k-R¹⁵ and
- (14) C₁₋₄alkyl, optionally substituted with from 1 to 6 halo groups,

R⁹ is selected from the group consisting of: hydrogen, C₁₋₁₂alkyl and aryl, wherein C₁₋₁₂alkyl and aryl are optionally substituted from one up to the maximum number of substituents with halo;

each R¹¹, R¹² and R¹⁶ is independently selected from the group consisting of:

- (1) hydrogen,
- (2) halo,
- (3) C₁₋₆alkyl,
- (4) C₂₋₆alkenyl,
- (5) C₁₋₆alkoxy and
- (6) hydroxy,

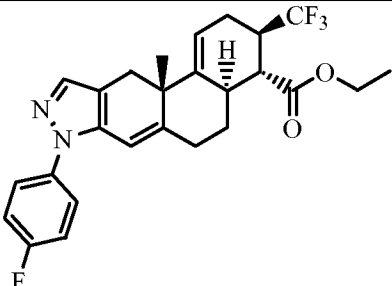
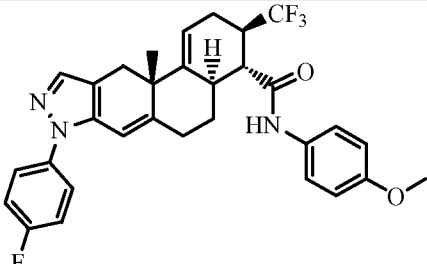
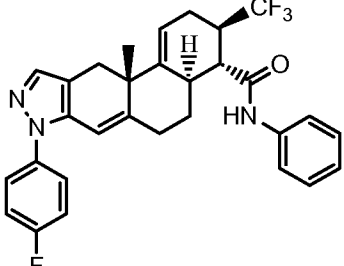
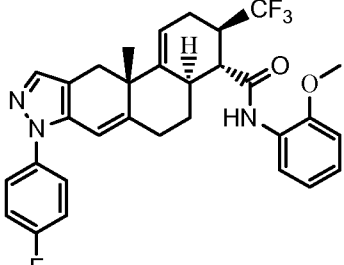
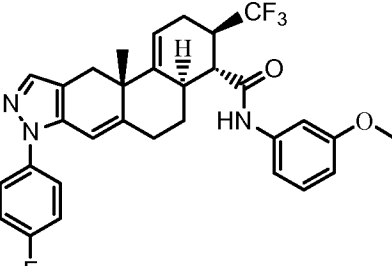
wherein items (3) to (5) above are optionally substituted from one up to the maximum number of substitutable positions with a substituent independently selected from the group consisting of: halo, OR¹², N(R¹³)₂ and C₁₋₆alkyl-S(O)_k-, wherein k is 0, 1 or 2;

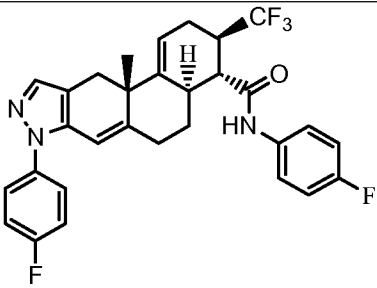
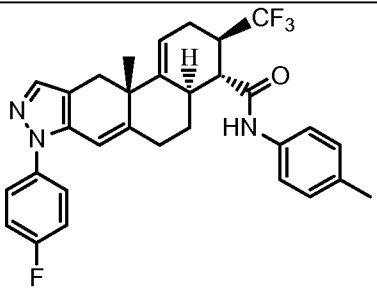
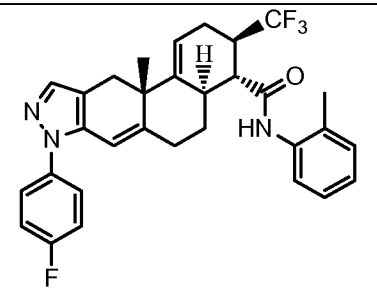
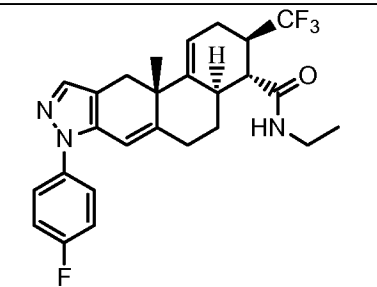
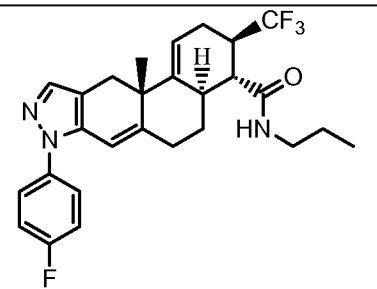
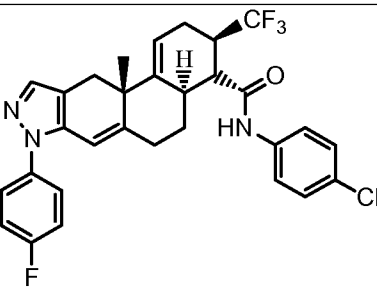
each R¹³ and R¹⁴ is independently selected from the group consisting of hydrogen and C₁₋₄alkyl, optionally substituted from one up to the maximum number of substitutable positions with halo; and

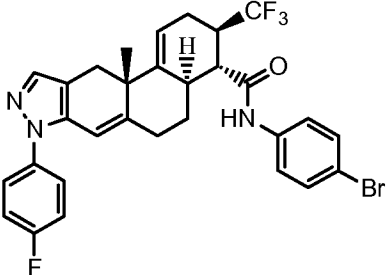
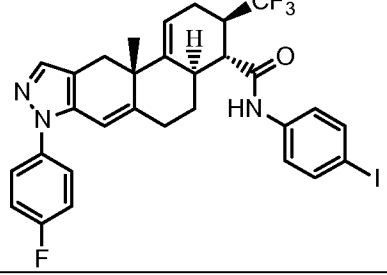
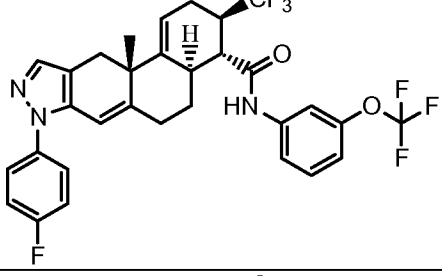
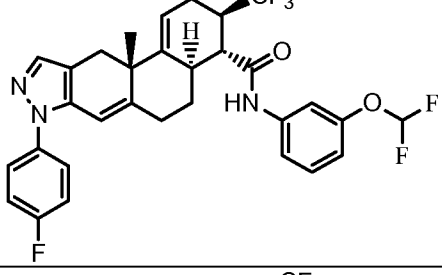
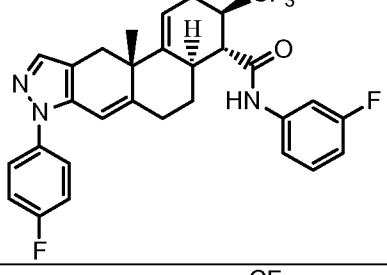
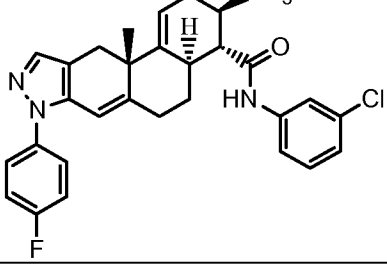
each R¹⁵ is independently selected from the group consisting of: hydrogen, C₁₋₆alkyl, aryl and C₁₋₁₂alkoxycarbonyl, wherein said C₁₋₆alkyl and C₁₋₁₂alkoxycarbonyl are optionally substituted from one up to the maximum number of substitutable positions with halo and said aryl is optionally substituted from one up to the maximum number of substitutable positions with halo and C₁₋₄alkyl, optionally substituted with 1-3 halo groups.

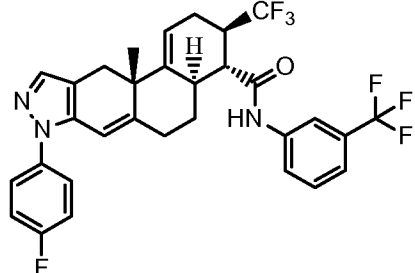
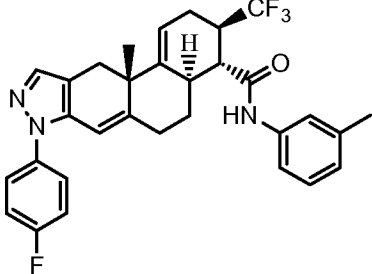
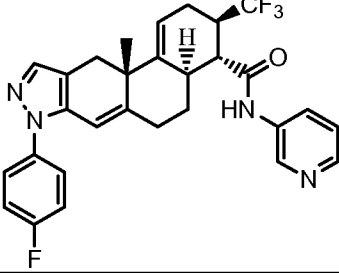
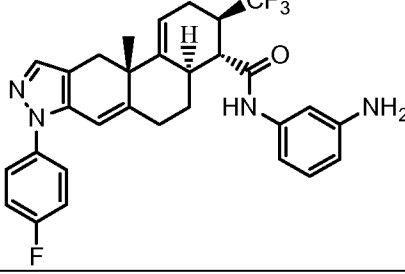
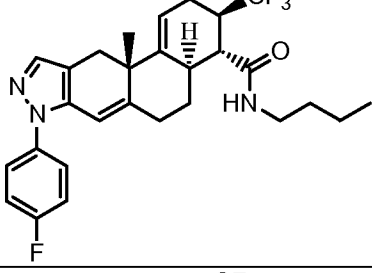
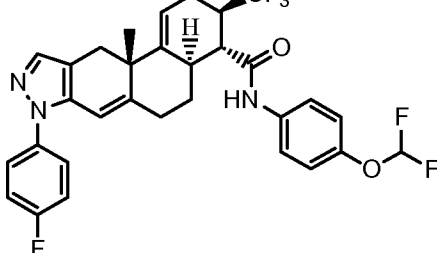
21. (Original) The compound according to Claim 20 wherein HET is selected from oxazolyl and imidazolyl.

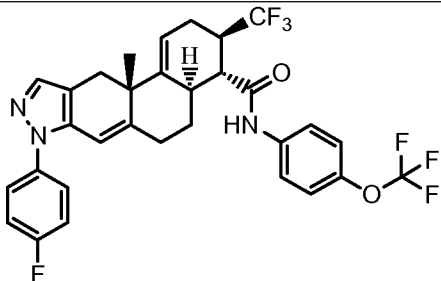
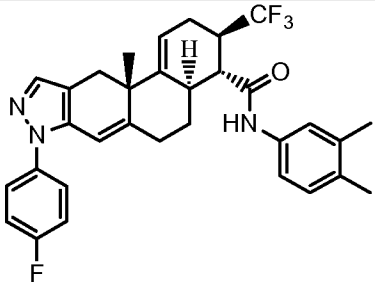
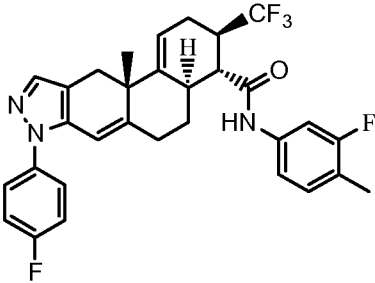
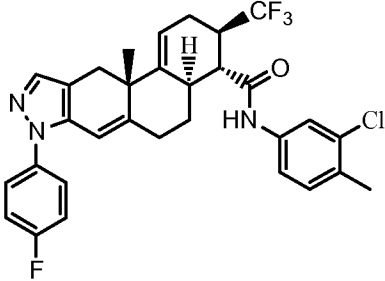
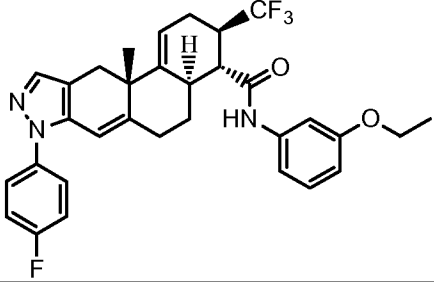
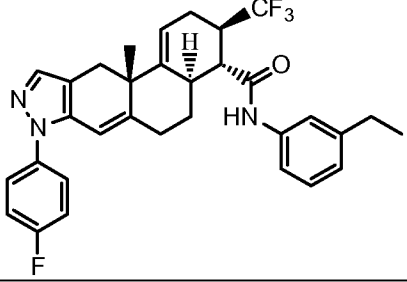
22. (Original) A compound selected from the group consisting of:

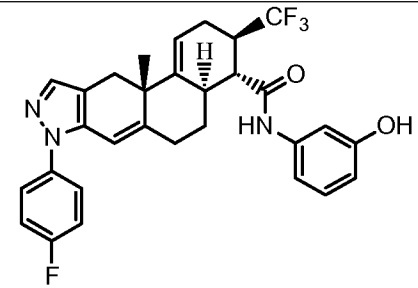
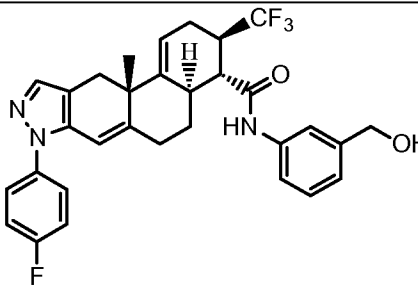
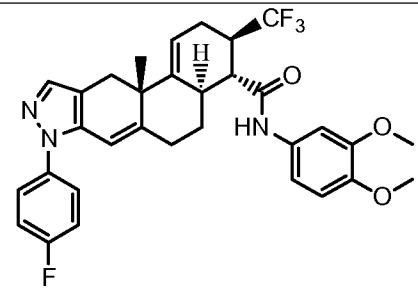
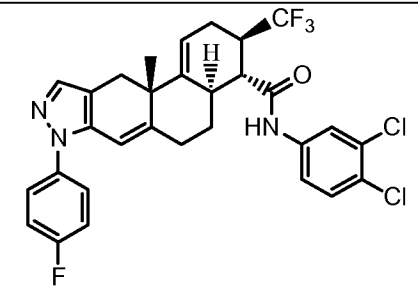
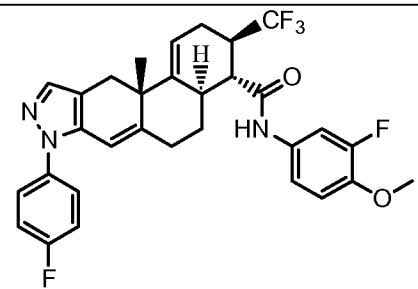
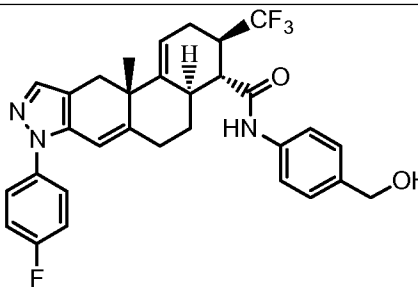
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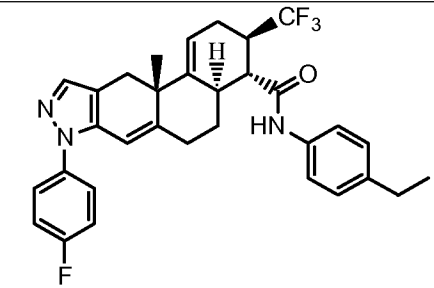
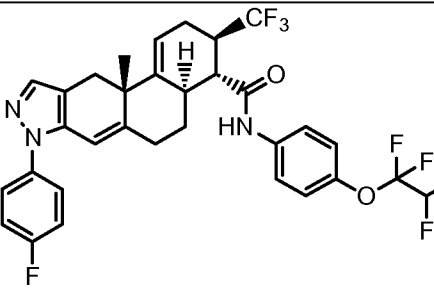
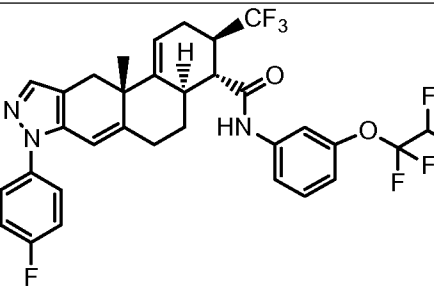
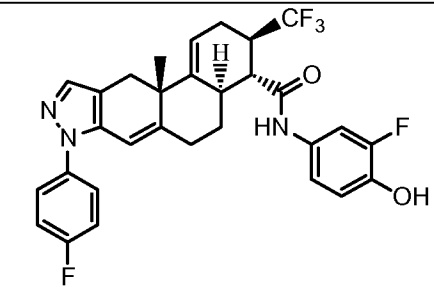
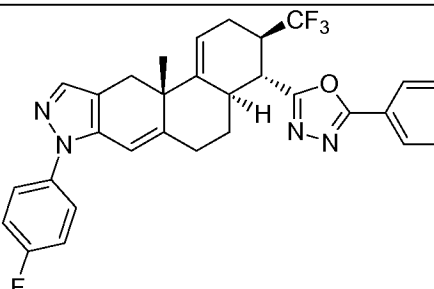
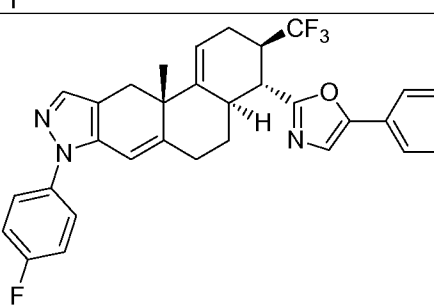
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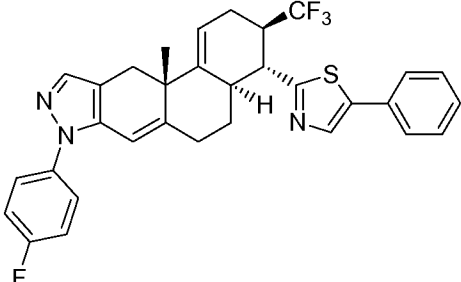
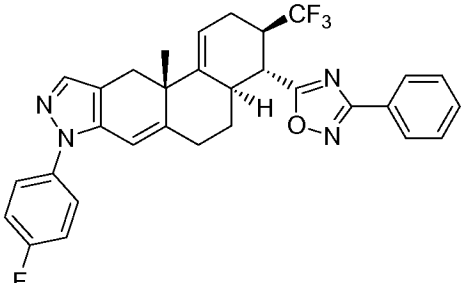
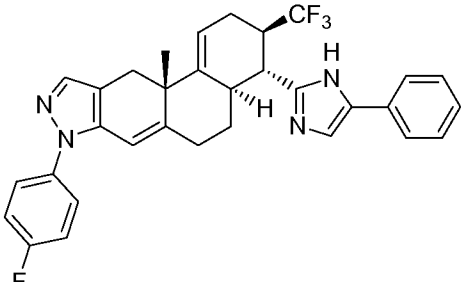
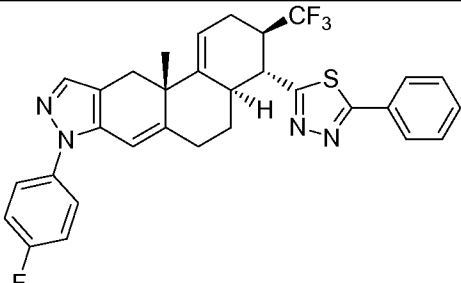
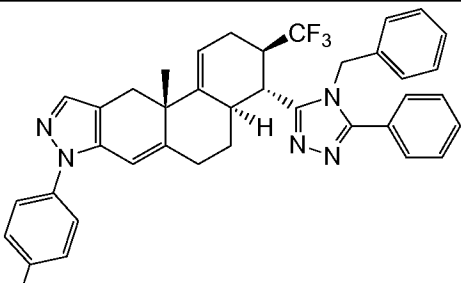
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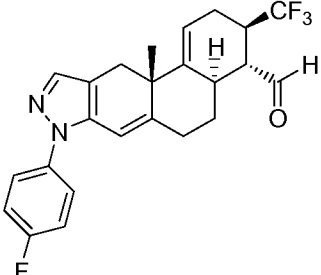
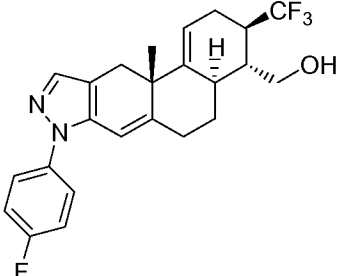
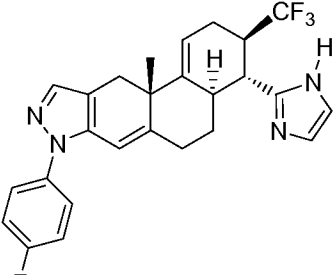
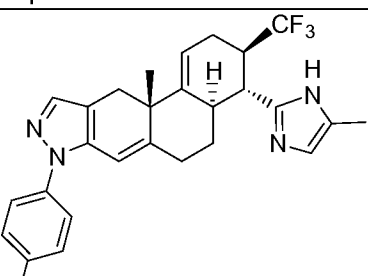
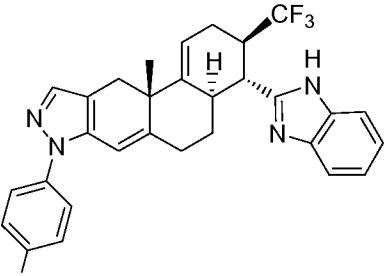
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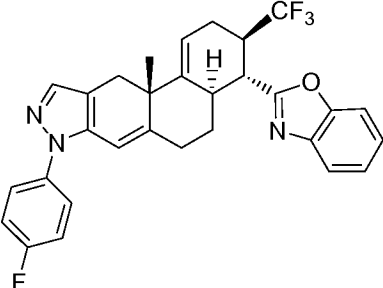
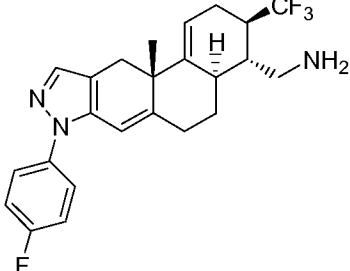
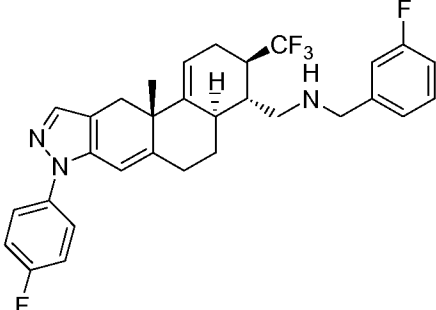
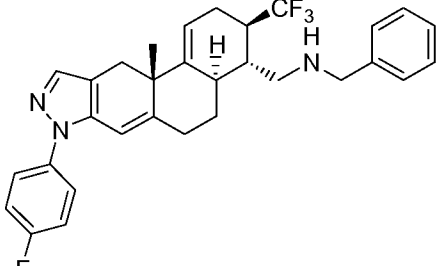
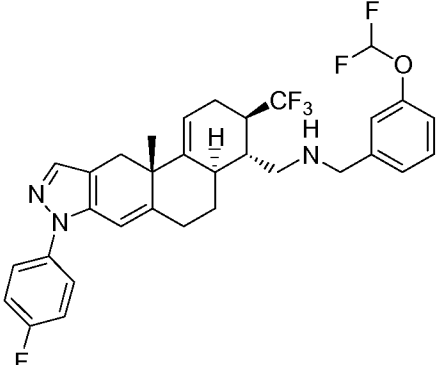
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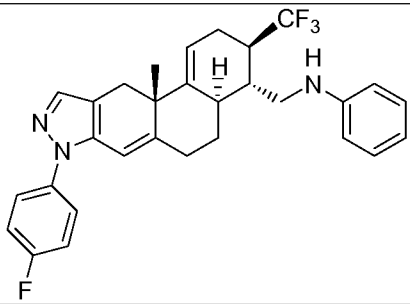
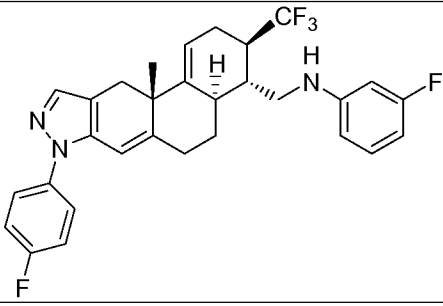
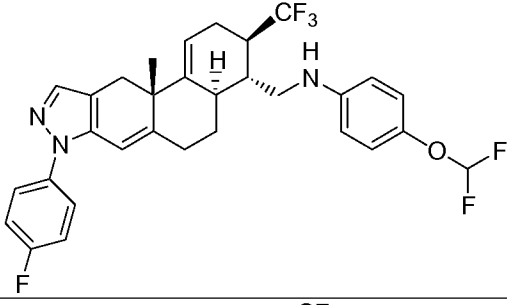
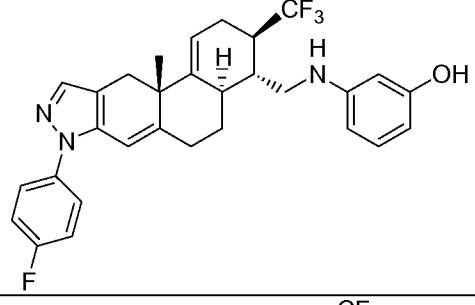
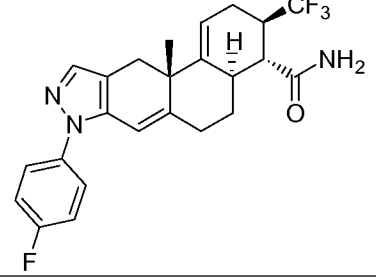
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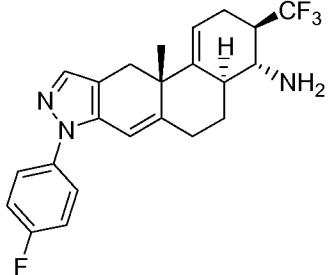
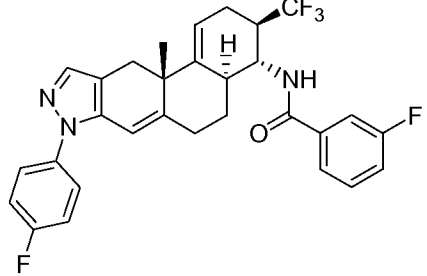
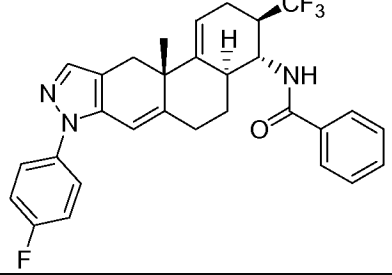
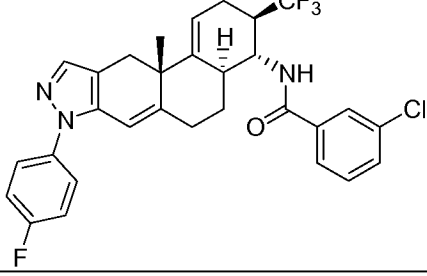
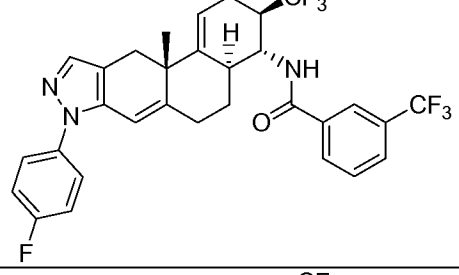
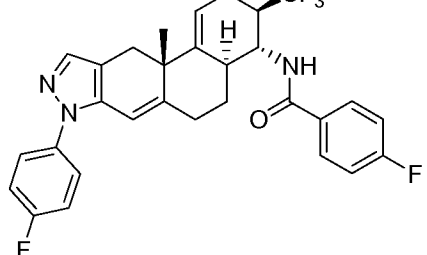
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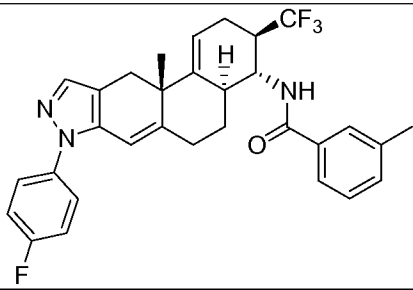
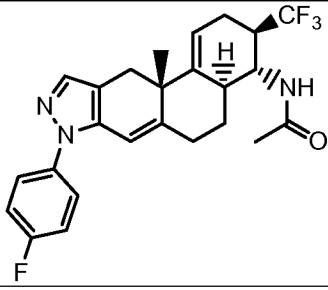
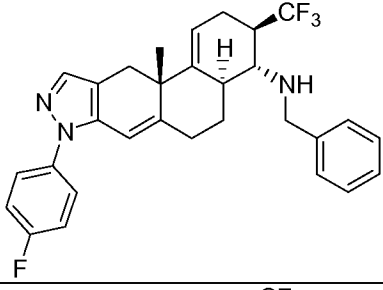
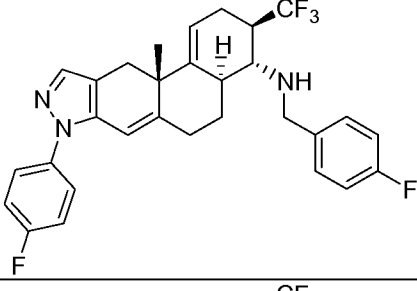
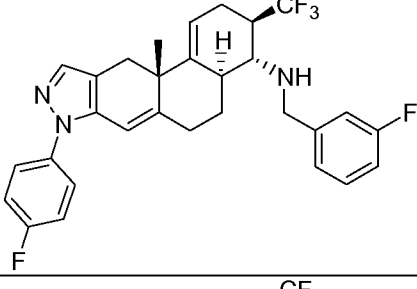
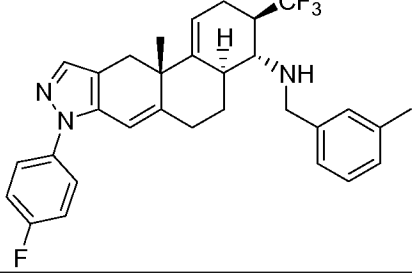
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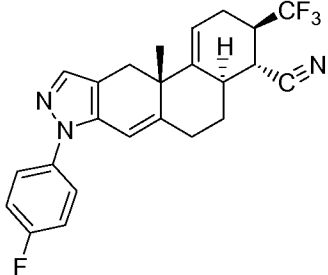
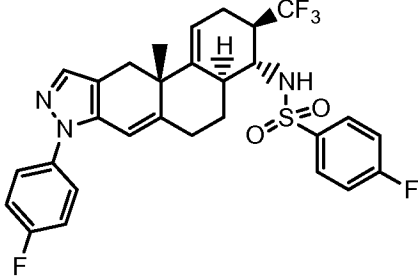
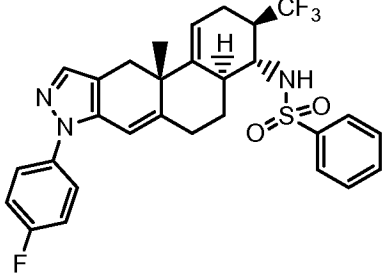
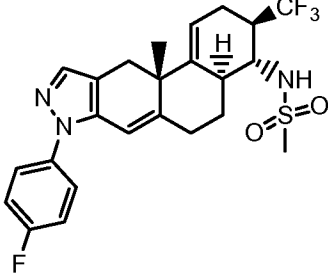
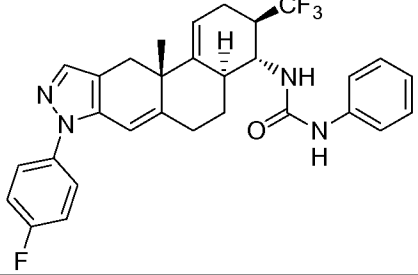
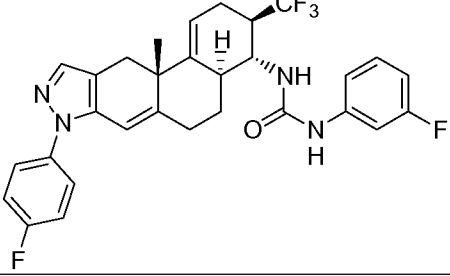
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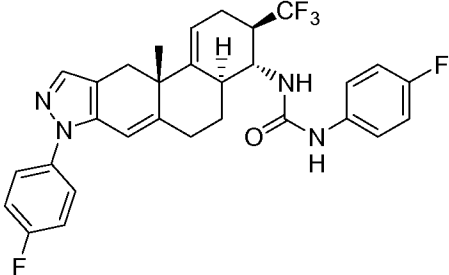
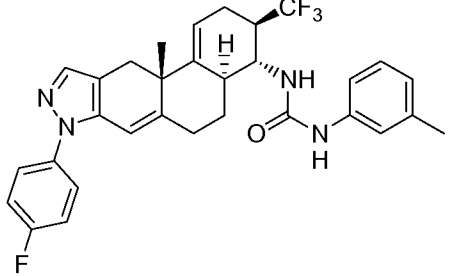
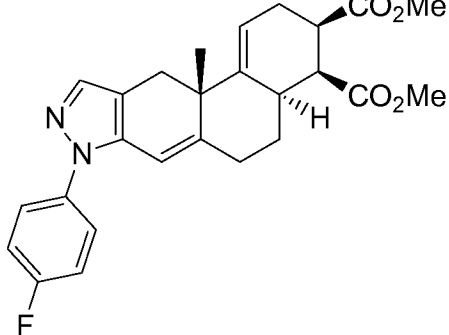
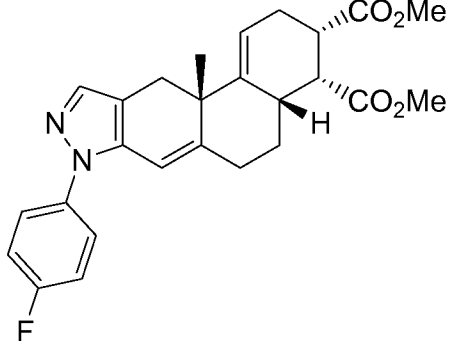
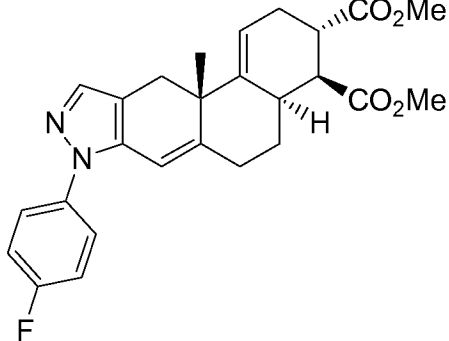
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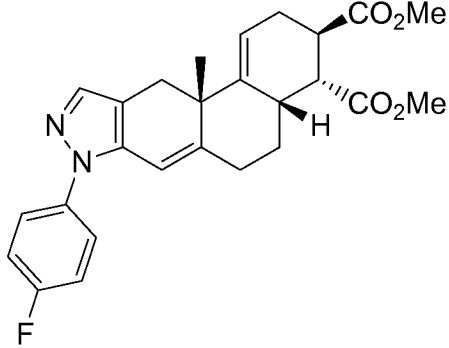
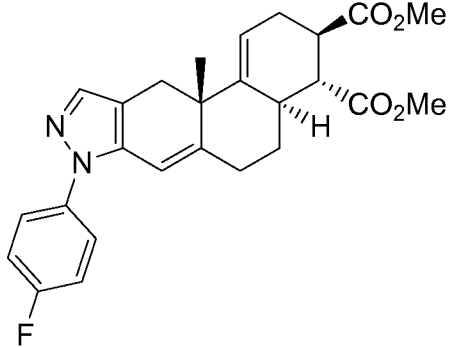
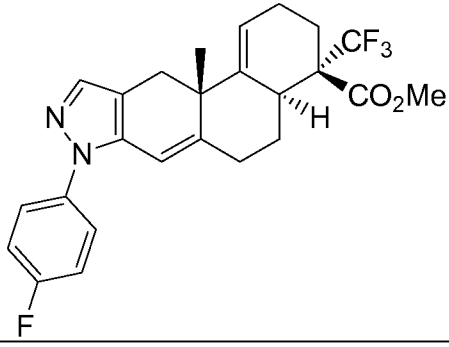
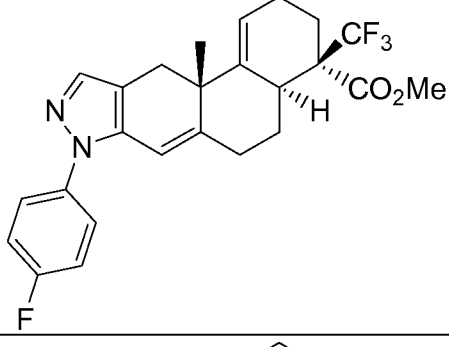
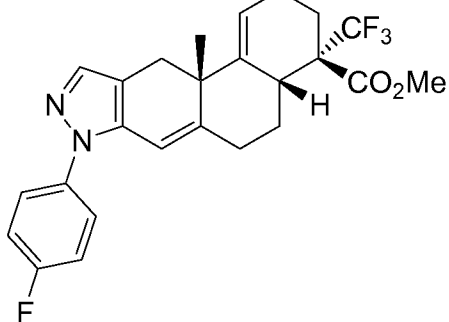
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58	 <chem>Fc1ccc(cc1)N2C3=C(C(=C2)C4C(C3)CC[C@H]4C)C(=O)N(Cc1ccc(F)cc1)C</chem>
59	 <chem>Fc1ccc(cc1)N2C3=C(C(=C2)C4C(C3)CC[C@H]4C)C(=O)N(Cc1ccc(OC(F)F)cc1)C</chem>
60	 <chem>Fc1ccc(cc1)N2C3=C(C(=C2)C4C(C3)CC[C@H]4C)C(=O)N(Cc1ccc(O)cc1)C</chem>
61	 <chem>Fc1ccc(cc1)N2C3=C(C(=C2)C4C(C3)CC[C@H]4C)C(=O)N(C)C</chem>

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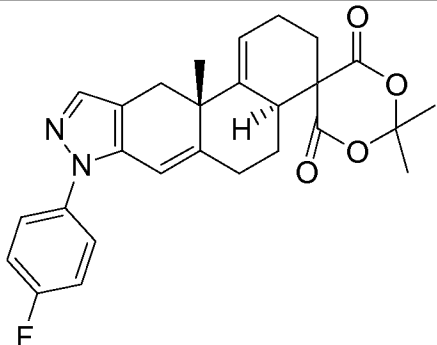
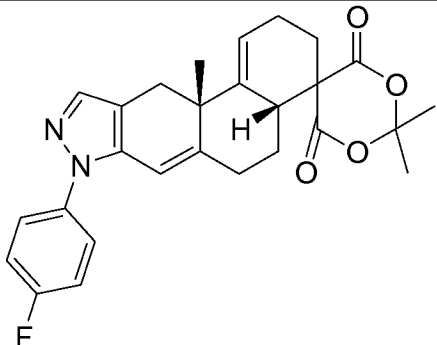
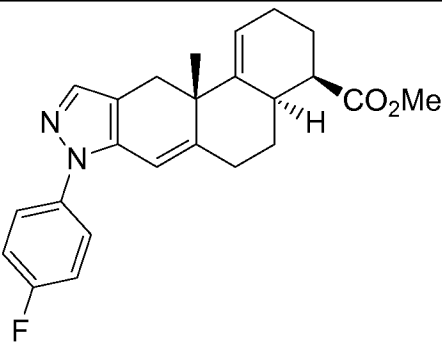
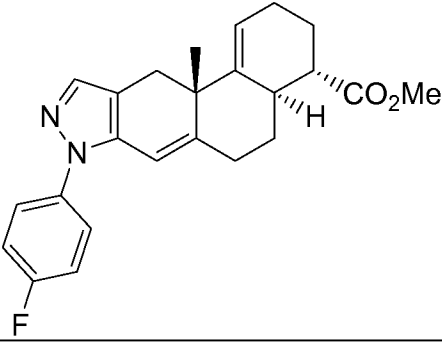
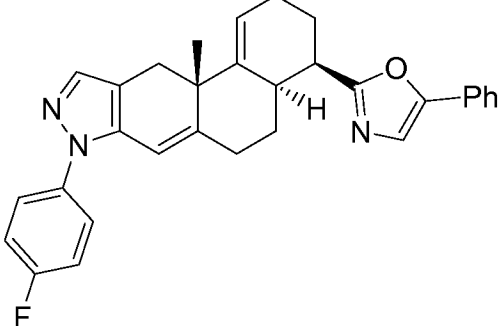
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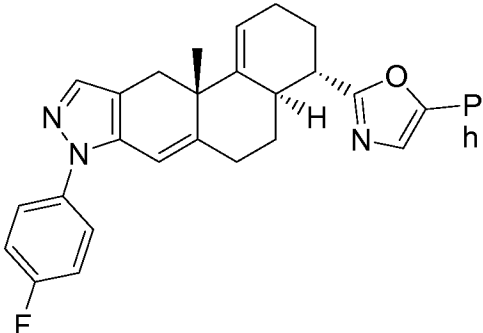
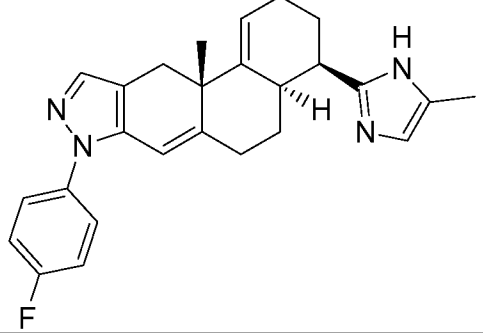
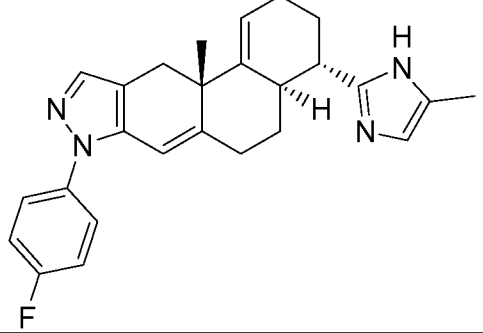
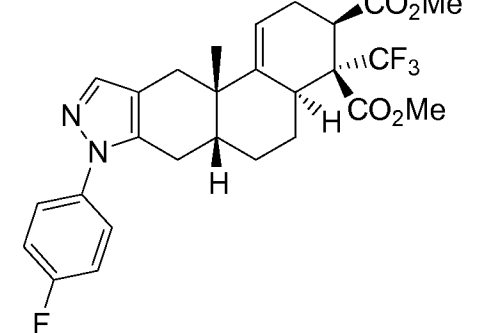
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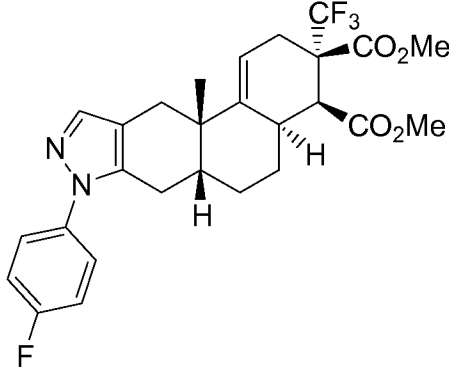
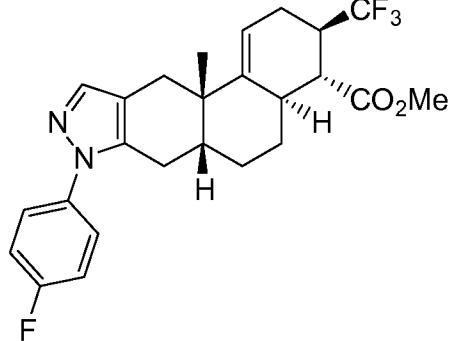
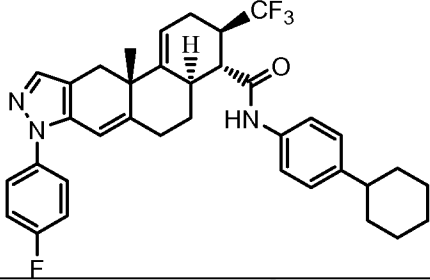
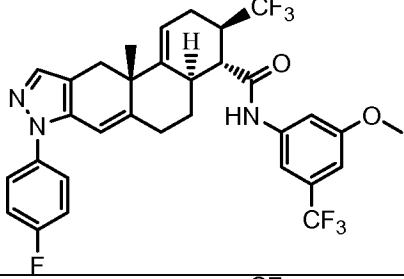
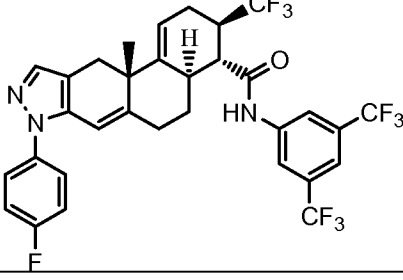
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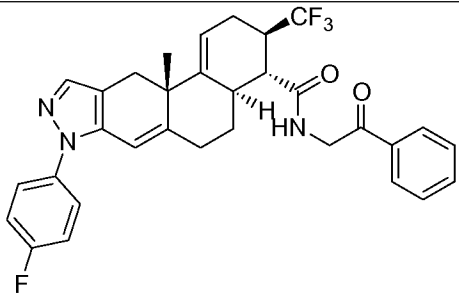
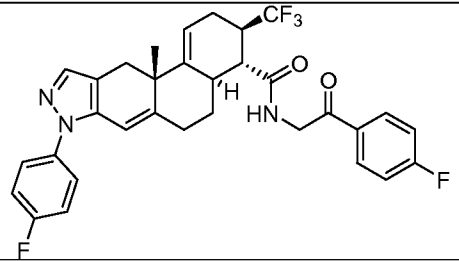
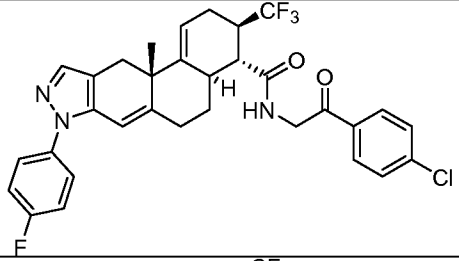
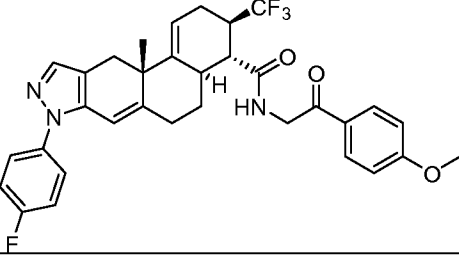
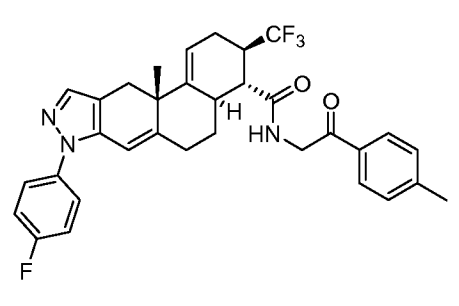
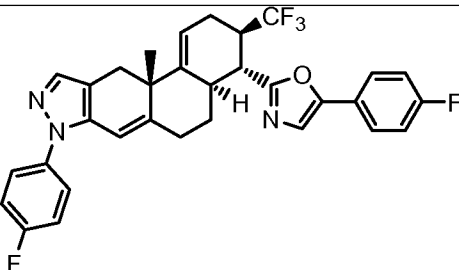
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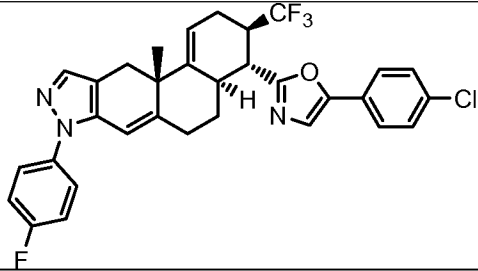
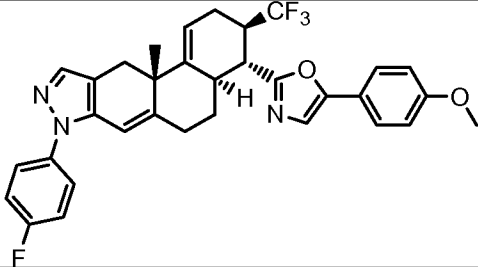
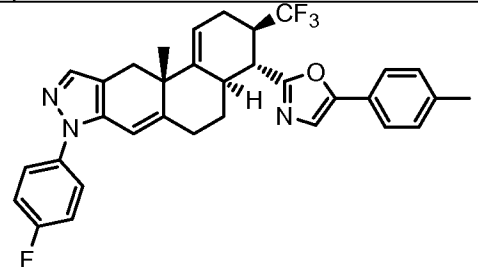
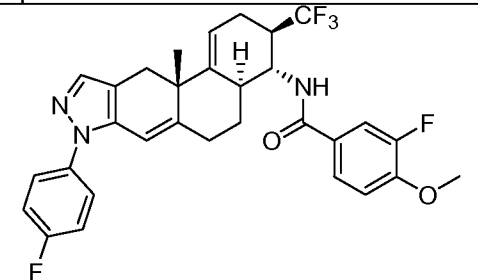
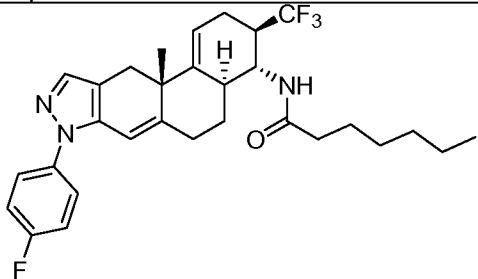
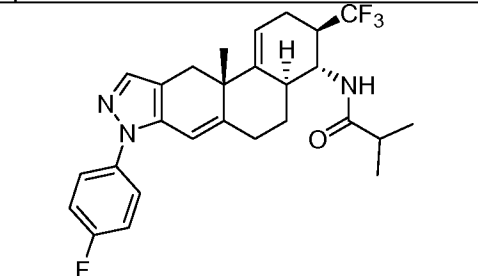
90	<p>Chemical structure 90: A steroid derivative. It features a 4-fluorophenylhydrazine group attached to the A-ring at C3. The B-ring has a double bond between C5 and C6. The D-ring has a methyl ester group (CO₂Me) at C17 and a trifluoromethyl group (CF₃) at C13. Stereochemistry is indicated with wedges and dashes.</p>
91	<p>Chemical structure 91: A steroid derivative. It features a 4-fluorophenylhydrazine group attached to the A-ring at C3. The B-ring has a double bond between C5 and C6. The D-ring has a methyl ester group (CO₂Me) at C17. Stereochemistry is indicated with wedges and dashes.</p>
92	<p>Chemical structure 92: A steroid derivative. It features a 4-fluorophenylhydrazine group attached to the A-ring at C3. The B-ring has a double bond between C5 and C6. The D-ring has a methyl ester group (CO₂Me) at C17. Stereochemistry is indicated with wedges and dashes.</p>
93	<p>Chemical structure 93: A steroid derivative. It features a 4-fluorophenylhydrazine group attached to the A-ring at C3. The B-ring has a double bond between C5 and C6. The D-ring has a methyl ester group (CO₂Me) at C17 and a trifluoromethyl group (CF₃) at C13. Stereochemistry is indicated with wedges and dashes.</p>
94	<p>Chemical structure 94: A steroid derivative. It features a 4-fluorophenylhydrazine group attached to the A-ring at C3. The B-ring has a double bond between C5 and C6. The D-ring has a methyl ester group (CO₂Me) at C17 and a trifluoromethyl group (CF₃) at C13. Stereochemistry is indicated with wedges and dashes.</p>

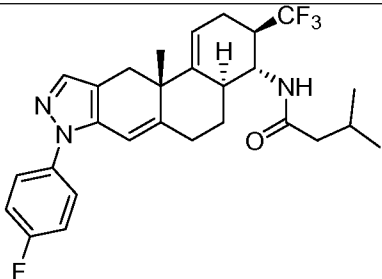
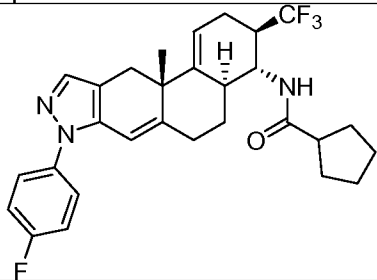
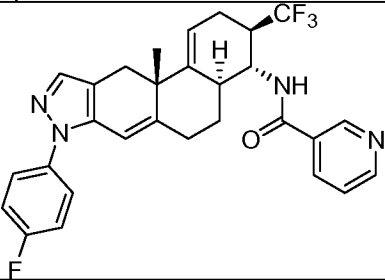
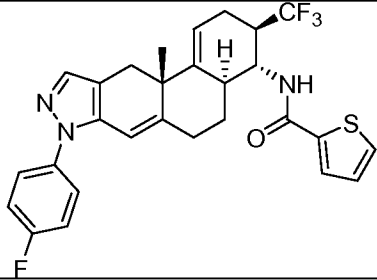
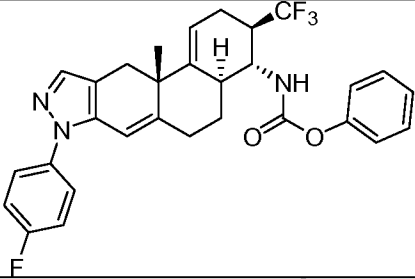
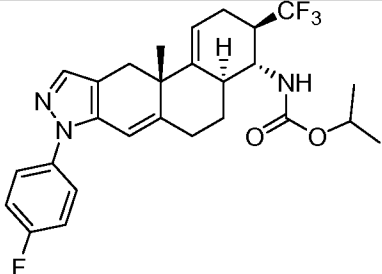
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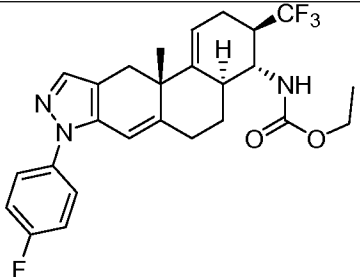
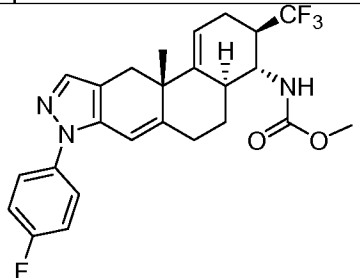
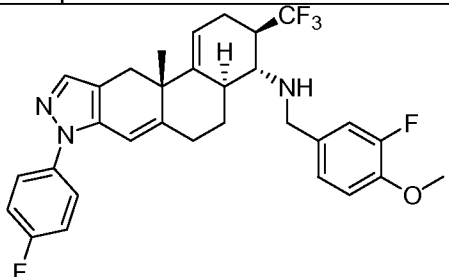
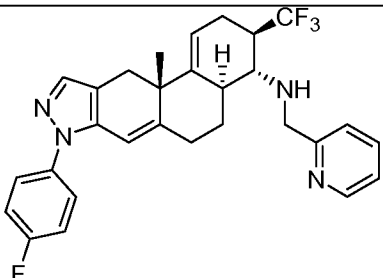
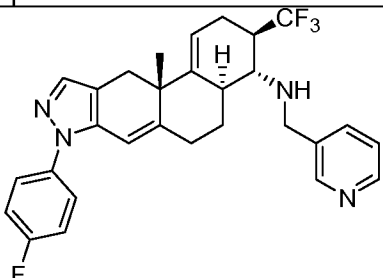
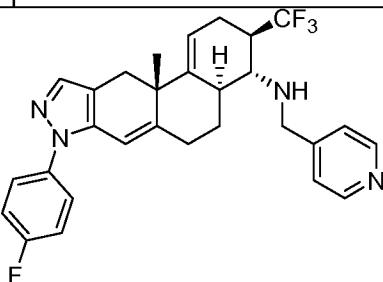
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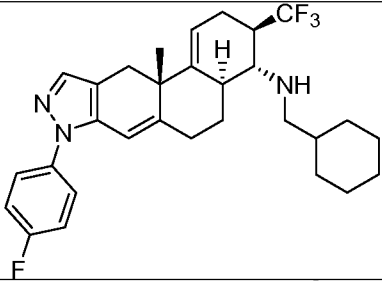
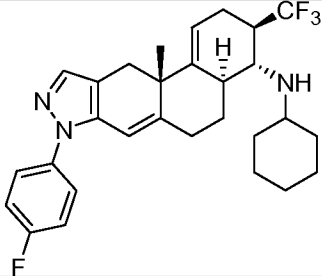
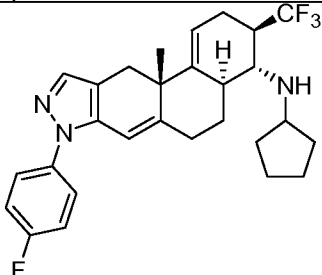
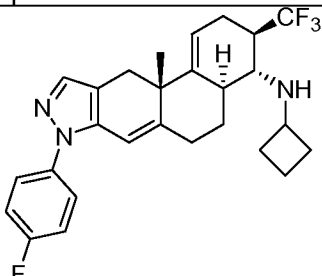
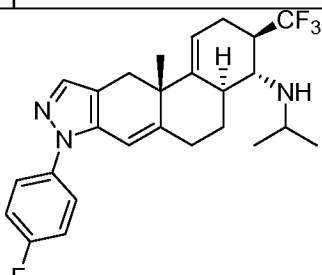
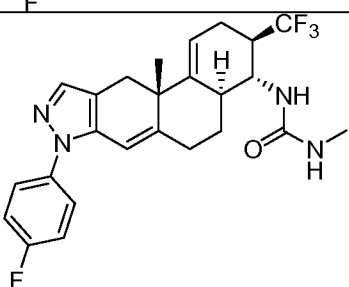
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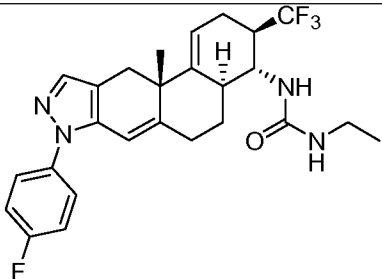
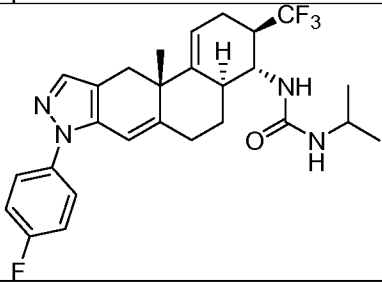
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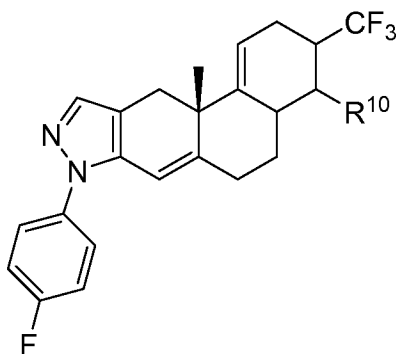
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23. (Original) A pharmaceutical composition comprising a compound according to Claim 1 in combination with a pharmaceutically acceptable carrier.

24. (Withdrawn) A method for treating a glucocorticoid receptor mediated disease or condition in a mammalian patient in need of such treatment comprising administering the patient a compound according to Claim 1 in an amount that is effective for treating the glucocorticoid receptor mediated disease or condition.

25-28. (Previously Canceled)

29. (Original) A compound according to Claim 1 of Formula Id



Id

or a pharmaceutically acceptable salt thereof, wherein

R¹⁰ is a 5-membered aromatic or non-aromatic mono-cyclic ring containing 1-3 heteroatoms selected from O, S, and N, and

R¹⁰ is mono-substituted with phenyl, wherein phenyl is optionally substituted with 1-3 substituents independently selected from halo, C₁₋₄alkyl and C₁₋₄alkoxy.

30. (Original) The compound according to Claim 29 wherein R¹⁰ is oxazolyl, oxadiazolyl or thiazolyl.

31. (Canceled)